

# Ketan S Khare

## List of Publications by Year in descending order

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Version: 2024-02-01

9  
papers

478  
citations

1306789

7  
h-index

1473754

9  
g-index

10  
all docs

10  
docs citations

10  
times ranked

630  
citing authors

#	ARTICLE	IF	CITATIONS
1	Atomistic simulation of volumetric properties of epoxy networks: effect of monomer length. <i>Soft Matter</i> , 2021, 17, 9957-9966.	1.2	4
2	Integration of Atomistic Simulation with Experiment Using Time~Temperature Superposition for a Cross~Linked Epoxy Network. <i>Macromolecular Theory and Simulations</i> , 2020, 29, 1900032.	0.6	12
3	Quantitative Comparison of Atomistic Simulations with Experiment for a Cross-Linked Epoxy: A Specific Volume~Cooling Rate Analysis. <i>Macromolecules</i> , 2018, 51, 564-575.	2.2	34
4	Bi-modal polymer networks: Composition-dependent trends in thermal, volumetric and structural properties from molecular dynamics simulation. <i>Polymer</i> , 2015, 58, 199-208.	1.8	23
5	Temperature dependence of creep compliance of highly cross-linked epoxy: A molecular simulation study. <i>AIP Conference Proceedings</i> , 2014, , .	0.3	2
6	Effect of Carbon Nanotube Functionalization on Mechanical and Thermal Properties of Cross-Linked Epoxy~Carbon Nanotube Nanocomposites: Role of Strengthening the Interfacial Interactions. <i>ACS Applied Materials &amp; Interfaces</i> , 2014, 6, 6098-6110.	4.0	163
7	High strain rate mechanical properties of a cross-linked epoxy across the glass transition. <i>Polymer</i> , 2013, 54, 7048-7057.	1.8	94
8	Effect of Carbon Nanotube Dispersion on Glass Transition in Cross-Linked Epoxy~Carbon Nanotube Nanocomposites: Role of Interfacial Interactions. <i>Journal of Physical Chemistry B</i> , 2013, 117, 7444-7454.	1.2	98
9	Directed Diffusion Approach for Preparing Atomistic Models of Crosslinked Epoxy for Use in Molecular Simulations. <i>Macromolecular Theory and Simulations</i> , 2012, 21, 322-327.	0.6	48